

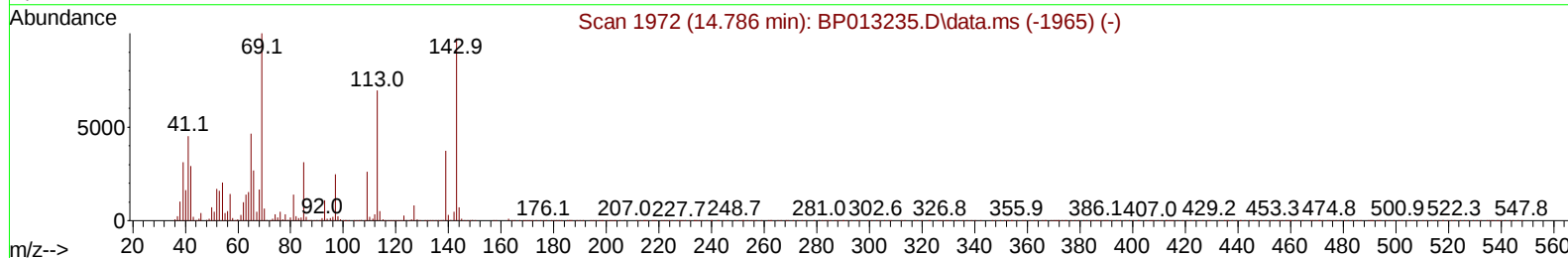
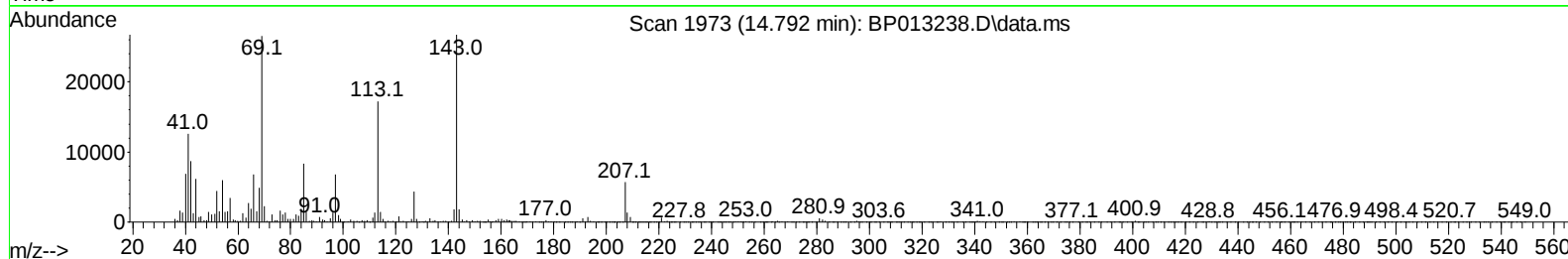
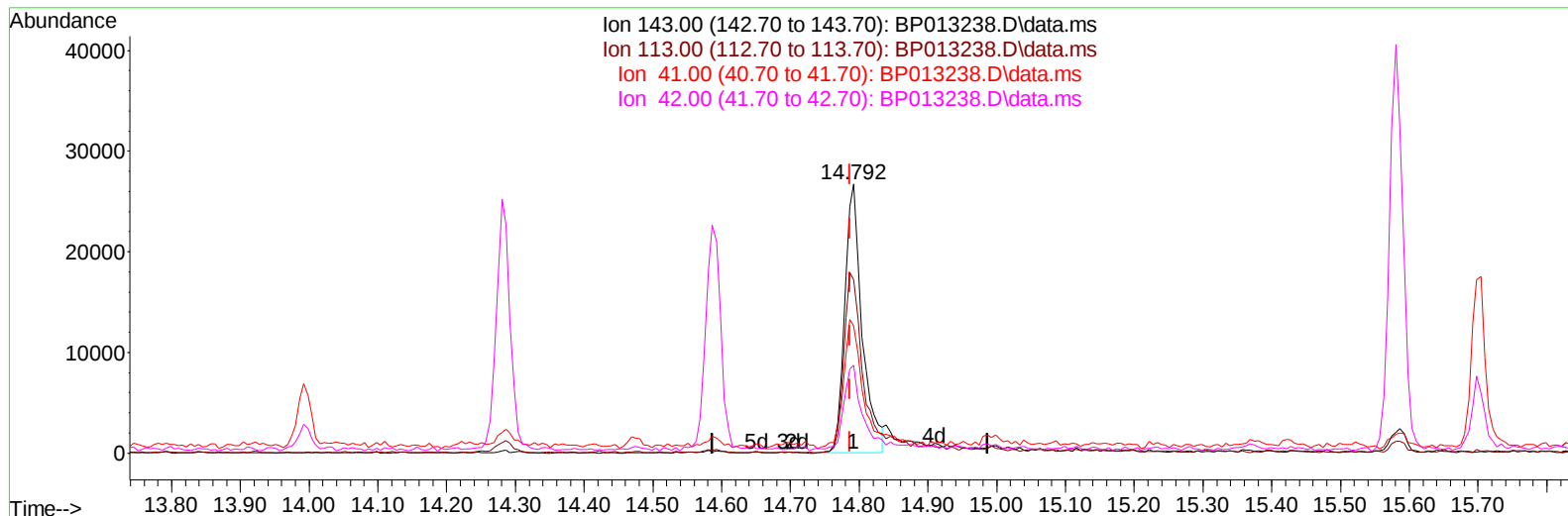
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013238.D
 Acq On : 22 Dec 2022 10:31
 Operator : CG/JU
 Sample : N6130-01 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXGJ0

Manual IntegrationsAPPROVED

Quant Time: Dec 22 21:44:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 23:38:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2022
 Supervised By :mohammad ahmed 12/23/2022



TIC: BP013238.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.006) 3.27 ng/u1

response 47287

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.80	64.56
41.00	35.70	47.21#
42.00	24.70	32.65#

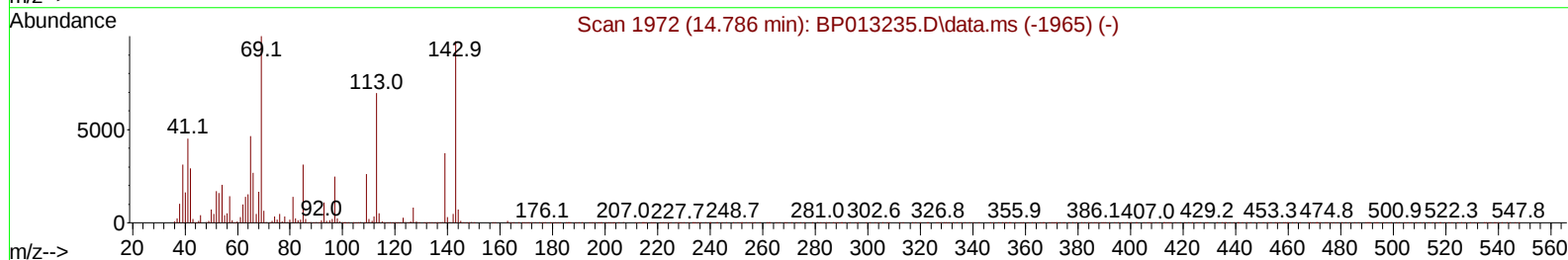
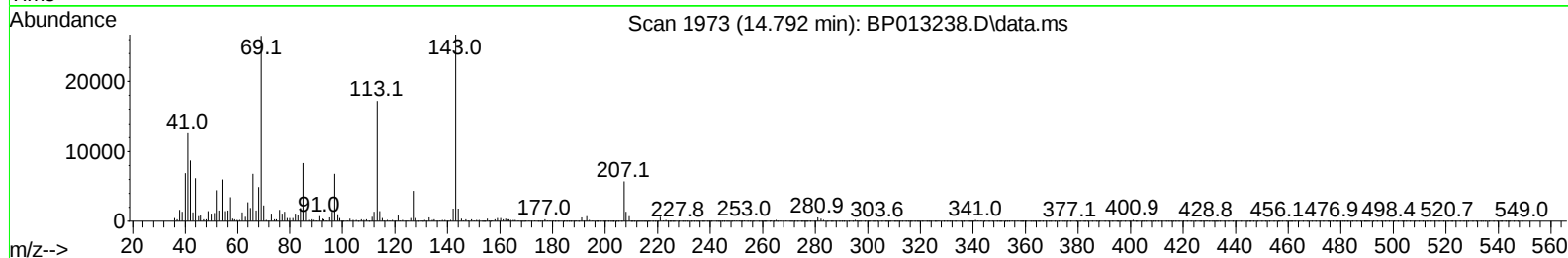
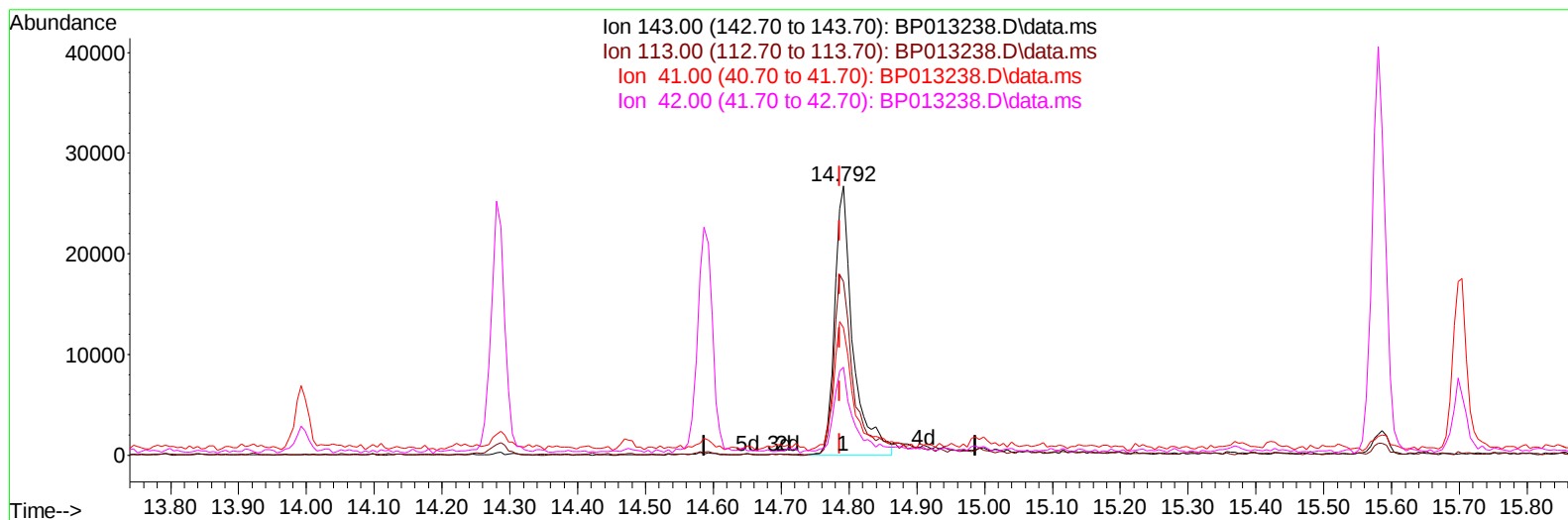
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TIC: BP013238.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.006) 3.49 ng/ul m

response 50430

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.80	64.56
41.00	35.70	47.21#
42.00	24.70	32.65#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	473341	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	1896098	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	1045288	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1968762	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	1975514	20.000	ng/ul	0.00
88) Perylene-d12	23.909	264	2286042	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	24942	2.253	ng/uL	0.00
4) Pyridine-d5	3.769	84	280275	8.912	ng/ul	0.00
7) Phenol-d5	7.104	99	247243	6.413	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	729742	31.376	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	621615	20.481	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	413633	13.093	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	423665	30.411	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	412466	26.299	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	712133	23.377	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	948565	22.056	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	2471177	30.761	ng/ul	0.00
49) Acenaphthylene-d8	14.286	160	2818394	31.927	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	50430m	3.490	ng/ul	0.00
60) Fluorene-d10	15.580	176	2168177	31.902	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	253606	20.017	ng/ul	0.00
73) Anthracene-d10	17.445	188	3446983	38.809	ng/ul	0.00
81) Pyrene-d10	19.680	212	4135015	36.466	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	4511717	39.595	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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